

A STUDY OF THE CRYSTAL STRUCTURE OF LATIFOLIN ( $C_{17}H_{18}O_4$ )



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#### ABSTRACT

X-Ray crystallographic data of Latifolin ( $C_{17}H_{18}O_4$ ) crystal were determined from rotation and Weissenberg photographs. The crystal belongs to the orthorhombic system with cell dimensions:

$$\begin{aligned}a &= 7.38 \pm 0.01 \text{ \AA}^{\circ} \\b &= 13.48 \pm 0.03 \text{ \AA}^{\circ} \\c &= 15.42 \pm 0.15 \text{ \AA}^{\circ}\end{aligned}$$

The space group is  $P2_1 2_1 2_1$ , the point group 2 2 2, there are four molecules per unit cell. The observed and calculated densities at  $27.5^{\circ} C$  are  $1.234 \text{ g/cm}^3$  and  $1.235 \text{ g/cm}^3$  respectively. Refined measurements of the unit cell dimension from a powder photograph yield the following values :

$$\begin{aligned}a &= 7.3887 \pm 0.0006 \text{ \AA}^{\circ} \\b &= 13.4581 \pm 0.0008 \text{ \AA}^{\circ} \\c &= 15.6157 \pm 0.0010 \text{ \AA}^{\circ}\end{aligned}$$

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## บทคัดย่อ

จากการใช้รังสีเอกซ์เคราะห์ของอุณหางคริสตอลโลกราฟฟี่ (Crystalllography) ของบล็อกแล็ทีโนเดิน (Latifolin) โดยใช้วิธีด้วยแบบการหมุนของผลึก (Rotation) และแบบไวอาเซ็นเบอร์ก (Weissenberg) พิจารณาผลลัพธ์ในระบบอธอรอมบิก (Orthorhombic system) มีค่าคงที่หน่วยเซลล์ (Cell dimensions)

|                      |           |
|----------------------|-----------|
| $a = 7.38 \pm 0.01$  | อัองสตรอน |
| $b = 13.48 \pm 0.03$ | อัองสตรอน |
| $c = 15.42 \pm 0.15$ | อัองสตรอน |

หมู่สี่มันนี้ ๓ มิติ (Space group) ของบล็อกเป็น  $P_{2_1} 2_1 2_1$  อยู่ในหมู่พอนท์ (Point group) ๒ ๒ ๒ และในหน่วยเซลล์มี ๔ โมเลกุล ความหนาแน่นของผลึกที่  $27.5^\circ C$  วัดได้  $1.234$  กรัม ต่อลูกบาศก์เซนติเมตร คำนวนได้  $1.235$  กรัม ต่อลูกบาศก์เซนติเมตร จากการหาค่าคงที่ของหน่วยเซลล์ให้ละเอียดขึ้น (Refined measurements) โดยใช้การถ่ายรูปแบบการกระจากรังสีเอกซ์วิวดาร์ทที่เป็นผง (Powder photograph) ได้

|                          |           |
|--------------------------|-----------|
| $a = 7.3887 \pm 0.0006$  | อัองสตรอน |
| $b = 13.4581 \pm 0.0008$ | อัองสตรอน |
| $c = 15.6157 \pm 0.0010$ | อัองสตรอน |

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